Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: SSSPTA1639MLS

PASSWORD:

l

TERMINAL (ENTER 1, 2, 3, OR ?):2

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NEWS 4 AUG 11 STN AnaVist workshops to be held in North America

NEWS 5 AUG 30 CA/CAplus -Increased access to 19th century research documents

NEWS 6 AUG 30 CASREACT - Enhanced with displayable reaction conditions

NEWS 7 SEP 09 ACD predicted properties enhanced in REGISTRY/ZREGISTRY

NEWS 8 SEP 22 MATHDI to be removed from STN

NEWS EXPRESS JUNE 13 CURRENT WINDOWS VERSION IS V8.0, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 13 JUNE 2005

NEWS HOURS STN Operating Hours Plus Help Desk Availability

NEWS INTER General Internet Information

NEWS LOGIN Welcome Banner and News Items

NEWS PHONE Direct Dial and Telecommunication Network Access to STN

NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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=> fil reg COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 0.21 0.21

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DICTIONARY FILE UPDATES: 25 SEP 2005 HIGHEST RN 863878-84-6

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Structure search iteration limits have been increased. See HELP SLIMITS for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> 437655-42-0/rn L2 1 437655-42-0/RN

=> d ibib abs 12
'IBIB' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'
'ABS' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN FIDE - All substance data, except seguence data

IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties
EPROP - Table of experimental properties

PROP - EPROP and CALC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

<code>HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats.</code>

ENTER DISPLAY FORMAT (IDE):bib, abs

'BIB' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

'ABS' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

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SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties EPROP - Table of experimental properties

PROP - EPROP and CALC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

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IPC -- International Patent Classification

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IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ---- AN, plus Bibliographic Data (original)

OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations

SIBIB ----- IBIB, no citations

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The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

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HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats.

ENTER DISPLAY FORMAT (IDE):bib

'BIB' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

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REG - RN

SAM - Index Name, MF, and structure - no RN

FIDE - All substance data, except sequence data

IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties

EPROP - Table of experimental properties

PROP - EPROP and CALC

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APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

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IABS -- ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original) OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats. ENTER DISPLAY FORMAT (IDE):sam

L2 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN

IN Acetamide, N,N'-(9,10-dihydro-9,10-dioxo-1,5-anthracenediyl)bis[2-mercapto-(9CI)

MF C18 H14 N2 O4 S2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

=> fil caplus COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 8.22 8.43

FULL ESTIMATED COST

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 12

L3 1 L2

=> d ibib abs 13

L3 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 2002:462494 CAPLUS

DOCUMENT NUMBER: 137:40117

TITLE: Tuned multifunctional linker molecules for electronic

charge transport through organic-inorganic composite

structures and use thereof

INVENTOR(S): Ford, William E.; Wessels, Jurina; Yasuda, Akio PATENT ASSIGNEE(S): Sony International (Europe) G.m.b.H., Germany

SOURCE: Eur. Pat. Appl., 36 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

	PATENT NO.				KIND DATE				APPLICATION NO.						DATE				
	EP	1215	 205	:		A1	-	2002	 0619		 EP 2	2000-	 1269	68		•	 20001	208	
		R:	•									•	LI,	LU,	NL,	SE	, MC,	PT,	
			ΙE,	SI,	LT,	LV,	FΙ,	RO,	MK,	CY,	AL,	TR							
	US	2002	1277	56		A1		2002	0912		US 2	2001-	6636				20011	206	
	JP	2002	2654	33		A2		2002	0918		JP 2	2001-	3749	16			20011	207	
PRIOR	RITY	APP	LN.	INFO	.:						EP 2	-000	1269	68		Α	20001	208	
AB	The	e pro	blem	und	erly:	ing t	the	pres	ent .	inve	ntic	n is	to	prov	ide	mul	tifun	ction	na.

AB The problem underlying the present invention is to provide multifunctional linker mols. for tuning the conductivity in nanoparticle-linker assemblies which

can be used in the formation of electronic networks and circuits and thin films of nanoparticles. The problem is solved according to the invention by providing a multifunctional linker mol. of the general structure CON1-FUNC1-X-FUNC2-CON2 in which X is the central body of the mol., FUNC1 and FUNC2 independently of each other are mol. groups introducing a dipole moment and/or capable of forming intermol. and/or intramol. H bonding networks, and CON1 and CON2 independently of each other are mol. groups binding to nanostructured units comprising metal and semiconductor materials.

REFERENCE COUNT: 12 THERE ARE 12 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> fil reg COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	3.10	11.53
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.73	-0.73

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Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> 437655-41-9/rn L4 1 437655-41-9/RN

=> sam
ENTER LOGIC EXPRESSION, QUERY NAME, OR (END):end
SEARCH ENDED BY USER

=> d sam 14

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN

IN Acetamide, N,N'-(9,10-dihydro-9,10-dioxo-1,4-anthracenediyl)bis[2-mercapto-(9CI)

MF C18 H14 N2 O4 S2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

=> fil caplus COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 2.20 13.73 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -0.73

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 14

L5 1 L4

=> 15 not 13

L6 0 L5 NOT L3

=> fil reg

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 0.45 14.18 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -0.73

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TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2005

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*

* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added, *
* effective March 20, 2005. A new display format, IDERL, is now *

* available and contains the CA role and document type information. *

Structure search iteration limits have been increased. See HELP SLIMITS for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> 433713-40-7/rn L7 1 433713-40-7/RN

=> fil caplus TOTAL SINCE FILE COST IN U.S. DOLLARS ENTRY SESSION 14.61 0.43 FULL ESTIMATED COST DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -0.73

FILE 'CAPLUS' ENTERED AT 15:41:33 ON 26 SEP 2005 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

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New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 17

=> 18 not 13

L9 1 L8 NOT L3

=> d ibib abs 19

L9 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 2002:462542 CAPLUS

DOCUMENT NUMBER: 137:15023

TITLE: Selective chemical sensors based on interlinked

nanoparticle assemblies

INVENTOR(S): Vossmeyer, Tobias; Besnard, Isabelle; Wessels, Jurina;

Ford, William; Yasuda, Akio

PATENT ASSIGNEE(S): Sony International (Europe) G.m.b.H., Germany

SOURCE: Eur. Pat. Appl., 37 pp.

CODEN: EPXXDW

DOCUMENT TYPE: LANGUAGE:

Patent English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

	PATENT NO.				KIND DATE			APPLICATION NO.						DATE				
	ĒΡ	1215	 485			A1	-	2002	0619	E	2	000-	1271	- - 49		2	0001	212
		R:	ΑT,	BE,	CH,	DE,	DK,	ES,	FR,	GB,	GR,	IT,	LI,	LU,	NL,	SE,	MC,	PT,
			ΙE,	SI,	LT,	LV,	FI,	RO,	MK,	CY, Z	λL,	TR						
	ΑU	2001	09708	33		A 5		2002	0613	A	J 2	001-	9708	3		2	0011	205
	ΑU	7811	66			B2		2005	0512									
	US	2002	13236	51		A1	:	2002	0919	U:	5 2	001-	1338	8 .		2	0011	211
	CN	1359	002			Α	:	2002	0717	CI	1 2	001-	1454	59		2	0011	212
	JР	2002	22861	L6		A2	:	2002	0814	J	2	001-	3791	39		2	0011	212
PRIOF	(TI	APP	LN.	NFO.	. :					E	2	000-	1271	49	P	A 2	0001	212
AB	The	inv	entic	on re	elate	es to	o a i	nanoj	parti	cle :	fil	m co	mpri:	sing	a na	nopa	arti	cle
	net	work	forr	ned o	of na	anopa	arti	cles	inte	erlin	ced	by	link	er mo	ols.	The	e li	nker
	mol	s. h	ave a	at le	east	two	lin	ker :	units	tha	c	an b	ind 1	to tl	he si	ırfa	ce o	f the
	nar	onar	ticle	9 5	By -	intro	duc	ina ·	selec	tivit	- 37-	enha	ncin	a un	its i	n +1	ne 1	inke

network formed of nanoparticles interlinked by linker mols. The linker mols. have at least two linker units that can bind to the surface of the nanoparticles. By introducing selectivity-enhancing units in the linker mol., the selectivity of the nanoparticle film towards target analytes can be enhanced. A fine-tuning of the selectivity can be achieved by including a fine-tuning unit in the vicinity of the selectivity-enhancing unit. The nanoparticle film can be used to produce chemical sensors which are selective and stable in their performance.

REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> fil reg				
COST IN U.S. DOLLARS	SINCE FILE	TOTAL		
·	ENTRY	SESSION		
FULL ESTIMATED COST	3.55	18.16		
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL		
	ENTRY	SESSION		
CA SUBSCRIBER PRICE	-0.73	-1.46		

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DICTIONARY FILE UPDATES: 25 SEP 2005 HIGHEST RN 863878-84-6

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*

* The CA roles and document type information have been removed from * the IDE default display format and the ED field has been added, * effective March 20, 2005. A new display format, IDERL, is now *

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Structure search iteration limits have been increased. See HELP SLIMITS for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> 437655-43-1/rn L10 1 437655-43-1/RN

=> fil caplus COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 0.43 18.59 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -1.46

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L11 1 L10

=> 111 not 13

L12 0 L11 NOT L3

=> fil reg

COST IN.U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 0.45 19.04

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION

CA SUBSCRIBER PRICE 0.00 -1.46

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STRUCTURE FILE UPDATES: 25 SEP 2005 HIGHEST RN 863878-84-6 DICTIONARY FILE UPDATES: 25 SEP 2005 HIGHEST RN 863878-84-6

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TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2005

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=> d sam 113

L13 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN

IN Carbamodithioic acid, 1,4-phenylenebis- (9CI)

MF C8 H8 N2 S4

CI COM

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

=> logoff y COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION		
FULL ESTIMATED COST	1.34	20.38		
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION		
CA SUBSCRIBER PRICE	0.00	-1.46		

STN INTERNATIONAL LOGOFF AT 15:44:47 ON 26 SEP 2005

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LOGINID: SSSPTA1639MLS

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

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NEWS "Ask CAS" for self-help around the clock 2

NEWS 3 JUL 20 Powerful new interactive analysis and visualization software, STN AnaVist, now available

NEWS AUG 11 STN AnaVist workshops to be held in North America

5 AUG 30 CA/CAplus -Increased access to 19th century research documents NEWS

NEWS 6 AUG 30 CASREACT - Enhanced with displayable reaction conditions

NEWS SEP 09 ACD predicted properties enhanced in REGISTRY/ZREGISTRY

SEP 22 MATHDI to be removed from STN NEWS

NEWS EXPRESS JUNE 13 CURRENT WINDOWS VERSION IS V8.0, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 13 JUNE 2005

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* * * * * * * * * * * * * * STN Columbus

FILE 'HOME' ENTERED AT 16:15:57 ON 26 SEP 2005

=> fil reg

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

0.21 0.21

FULL ESTIMATED COST

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Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> d sam 11

L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN IN Acetamide, N,N'-1,4-cyclohexanediylbis[2-mercapto- (9CI) MF C10 H18 N2 O2 S2

$$\begin{matrix} \text{O} \\ \parallel \\ \text{HS-CH}_2\text{-C-NH} \end{matrix}$$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

=> d sam 12

L2 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN IN Carbamodithioic acid, 1,4-cyclohexanediylbis- (9CI) MF C8 H14 N2 S4

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

. => logoff y
 COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 3.11 3.32

STN INTERNATIONAL LOGOFF AT 16:17:37 ON 26 SEP 2005

Welcome to STN International! Enter x:x

LOGINID: SSSPTA1639MLS

PASSWORD:

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NEWS 8 SEP 22 MATHDI to be removed from STN

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=> fil req

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FULL ESTIMATED COST 0.21 0.21

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem. STRUCTURE FILE UPDATES: 25 SEP 2005 HIGHEST RN 863878-84-6 DICTIONARY FILE UPDATES: 25 SEP 2005 HIGHEST RN 863878-84-6 New CAS Information Use Policies, enter HELP USAGETERMS for details. TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2005 Please note that search-term pricing does apply when conducting SmartSELECT searches. ******************* * The CA roles and document type information have been removed from * * the IDE default display format and the ED field has been added, * effective March 20, 2005. A new display format, IDERL, is now * available and contains the CA role and document type information. * *********** Structure search iteration limits have been increased. See HELP SLIMITS for details. Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html => 1(w) 4(w) dimercaptoacetamido(w) 9(w) 10(w) anthraquinone 17900944 1 15963747 4 0 DIMERCAPTOACETAMIDO 1954377 9 1338254 10 12330 ANTHRAQUINONE 0 1 (W) 4 (W) DIMERCAPTOACETAMIDO (W) 9 (W) 10 (W) ANTHRAQUINONE L1=> 1(w)4(w)dimercaptoacetamidocyclohexane 17900944 1 15963747 4 0 DIMERCAPTOACETAMIDOCYCLOHEXANE L2 0 1 (W) 4 (W) DIMERCAPTOACETAMIDOCYCLOHEXANE => 1(w)5(w)dimercaptoacetamido(w)9(w)10(w)anthraquinone 17900944 1 10527616 5 0 DIMERCAPTOACETAMIDO 1954377 9 1338254 10 12330 ANTHRAQUINONE 0 1 (W) 5 (W) DIMERCAPTOACETAMIDO (W) 9 (W) 10 (W) ANTHRAQUINONE L3 => 1(w)8(w)dimercaptoacetamidooctane 17900944 1 2538423 8 0 DIMERCAPTOACETAMIDOOCTANE

=> 1(w)4(w)dithiocarbamatocyclohexane 17900944 1

0 1 (W) 8 (W) DIMERCAPTOACETAMIDOOCTANE

L4

15963747 4

0 DITHIOCARBAMATOCYCLOHEXANE

L5 0 1 (W) 4 (W) DITHIOCARBAMATOCYCLOHEXANE

=> dithiocarbamatocyclohexane

L6 0 DITHIOCARBAMATOCYCLOHEXANE

=> dimercaptoacetamidooctane

L7 0 DIMERCAPTOACETAMIDOOCTANE

=> dimercaptoacetamido

L8 0 DIMERCAPTOACETAMIDO

=> logoff y

COST IN U.S. DOLLARS

SINCE FILE TOTAL

ENTRY SESSION 113.84 114.05

FULL ESTIMATED COST

STN INTERNATIONAL LOGOFF AT 14:05:14 ON 26 SEP 2005